



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:57 PM UTC

PDB ID : 5NF1 / pdb_00005nf1
Title : crystal structure of variants
Authors : Linde, M.; Rajendran, C.; Babinger, P.; Sterner, R.
Deposited on : 2017-03-13
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

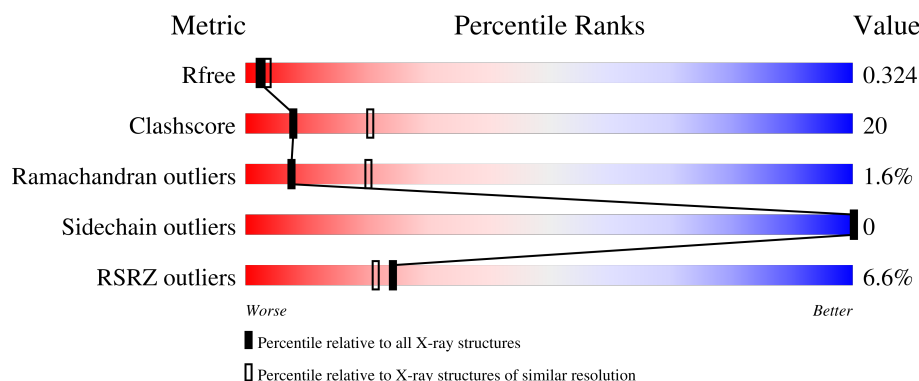
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	256	3% 56% 34% 9%
1	B	256	5% 54% 32% 14%
1	C	256	3% 65% 22% 12%
1	D	256	6% 56% 31% 12%
1	E	256	11% 53% 39% 7%

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Mol	Chain	Length	Quality of chain
1	F	256	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '7%', followed by a green segment labeled '54%', a yellow segment labeled '37%', and a small grey segment at the end labeled '10%'.

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9975 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Geranylgeranylglyceryl phosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1730	1089	294	337	10			
1	B	221	Total	C	N	O	S	0	0	0
			1590	1002	266	314	8			
1	C	226	Total	C	N	O	S	0	0	0
			1628	1036	273	309	10			
1	D	224	Total	C	N	O	S	0	0	0
			1610	1017	276	308	9			
1	E	237	Total	C	N	O	S	0	0	0
			1708	1073	286	339	10			
1	F	231	Total	C	N	O	S	0	0	0
			1666	1055	277	325	9			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP O26652
A	2	PHE	-	expression tag	UNP O26652
A	3	LYS	-	expression tag	UNP O26652
A	141	ALA	TRP	conflict	UNP O26652
A	249	LEU	-	expression tag	UNP O26652
A	250	GLU	-	expression tag	UNP O26652
A	251	HIS	-	expression tag	UNP O26652
A	252	HIS	-	expression tag	UNP O26652
A	253	HIS	-	expression tag	UNP O26652
A	254	HIS	-	expression tag	UNP O26652
A	255	HIS	-	expression tag	UNP O26652
A	256	HIS	-	expression tag	UNP O26652
B	1	MET	-	initiating methionine	UNP O26652
B	2	PHE	-	expression tag	UNP O26652
B	3	LYS	-	expression tag	UNP O26652
B	141	ALA	TRP	conflict	UNP O26652
B	249	LEU	-	expression tag	UNP O26652

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Chain	Residue	Modelled	Actual	Comment	Reference
B	250	GLU	-	expression tag	UNP O26652
B	251	HIS	-	expression tag	UNP O26652
B	252	HIS	-	expression tag	UNP O26652
B	253	HIS	-	expression tag	UNP O26652
B	254	HIS	-	expression tag	UNP O26652
B	255	HIS	-	expression tag	UNP O26652
B	256	HIS	-	expression tag	UNP O26652
C	1	MET	-	initiating methionine	UNP O26652
C	2	PHE	-	expression tag	UNP O26652
C	3	LYS	-	expression tag	UNP O26652
C	141	ALA	TRP	conflict	UNP O26652
C	249	LEU	-	expression tag	UNP O26652
C	250	GLU	-	expression tag	UNP O26652
C	251	HIS	-	expression tag	UNP O26652
C	252	HIS	-	expression tag	UNP O26652
C	253	HIS	-	expression tag	UNP O26652
C	254	HIS	-	expression tag	UNP O26652
C	255	HIS	-	expression tag	UNP O26652
C	256	HIS	-	expression tag	UNP O26652
D	1	MET	-	initiating methionine	UNP O26652
D	2	PHE	-	expression tag	UNP O26652
D	3	LYS	-	expression tag	UNP O26652
D	141	ALA	TRP	conflict	UNP O26652
D	249	LEU	-	expression tag	UNP O26652
D	250	GLU	-	expression tag	UNP O26652
D	251	HIS	-	expression tag	UNP O26652
D	252	HIS	-	expression tag	UNP O26652
D	253	HIS	-	expression tag	UNP O26652
D	254	HIS	-	expression tag	UNP O26652
D	255	HIS	-	expression tag	UNP O26652
D	256	HIS	-	expression tag	UNP O26652
E	1	MET	-	initiating methionine	UNP O26652
E	2	PHE	-	expression tag	UNP O26652
E	3	LYS	-	expression tag	UNP O26652
E	141	ALA	TRP	conflict	UNP O26652
E	249	LEU	-	expression tag	UNP O26652
E	250	GLU	-	expression tag	UNP O26652
E	251	HIS	-	expression tag	UNP O26652
E	252	HIS	-	expression tag	UNP O26652
E	253	HIS	-	expression tag	UNP O26652
E	254	HIS	-	expression tag	UNP O26652
E	255	HIS	-	expression tag	UNP O26652

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Chain	Residue	Modelled	Actual	Comment	Reference
E	256	HIS	-	expression tag	UNP O26652
F	1	MET	-	initiating methionine	UNP O26652
F	2	PHE	-	expression tag	UNP O26652
F	3	LYS	-	expression tag	UNP O26652
F	141	ALA	TRP	conflict	UNP O26652
F	249	LEU	-	expression tag	UNP O26652
F	250	GLU	-	expression tag	UNP O26652
F	251	HIS	-	expression tag	UNP O26652
F	252	HIS	-	expression tag	UNP O26652
F	253	HIS	-	expression tag	UNP O26652
F	254	HIS	-	expression tag	UNP O26652
F	255	HIS	-	expression tag	UNP O26652
F	256	HIS	-	expression tag	UNP O26652

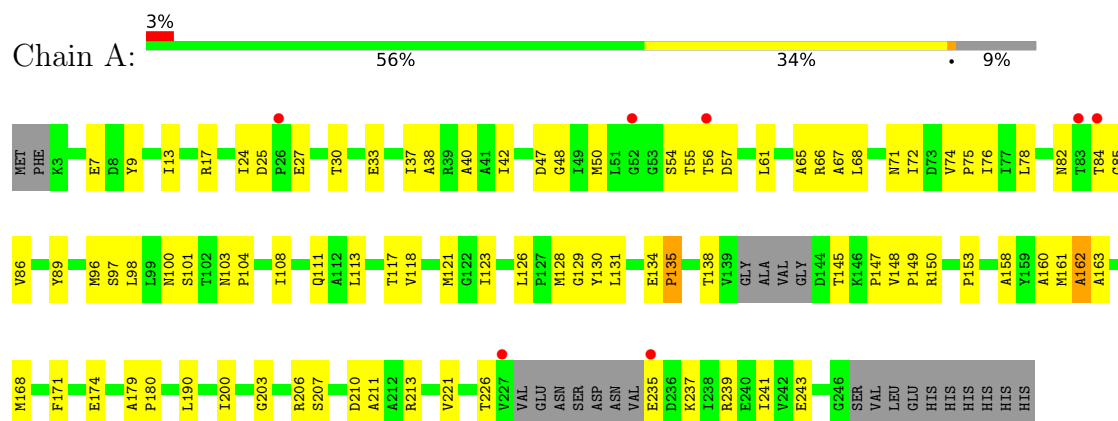
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	8	Total O 8 8	0	0
2	B	6	Total O 6 6	0	0
2	C	20	Total O 20 20	0	0
2	D	5	Total O 5 5	0	0
2	E	2	Total O 2 2	0	0
2	F	2	Total O 2 2	0	0

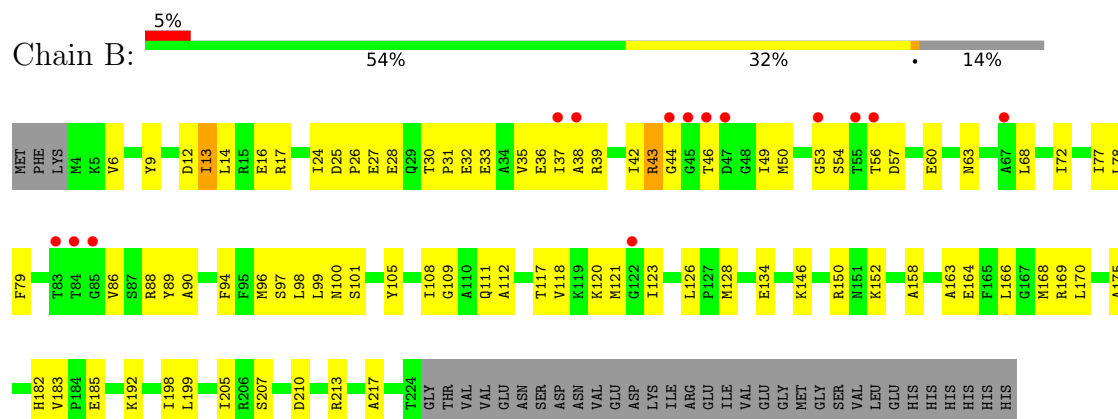
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

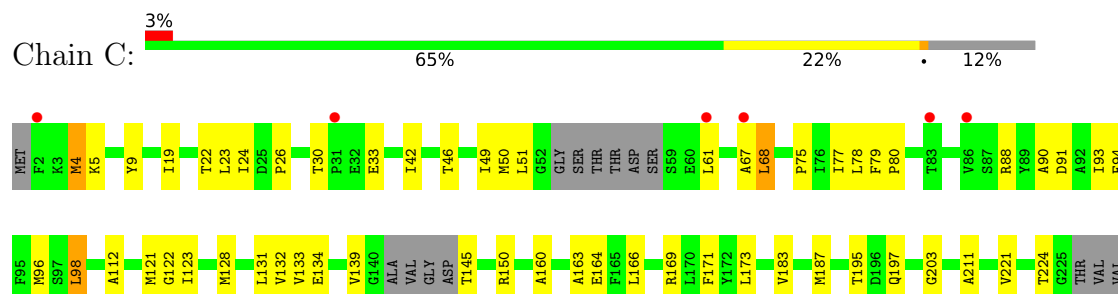
• Molecule 1: Geranylgeranylglyceryl phosphate synthase



• Molecule 1: Geranylgeranylglyceryl phosphate synthase

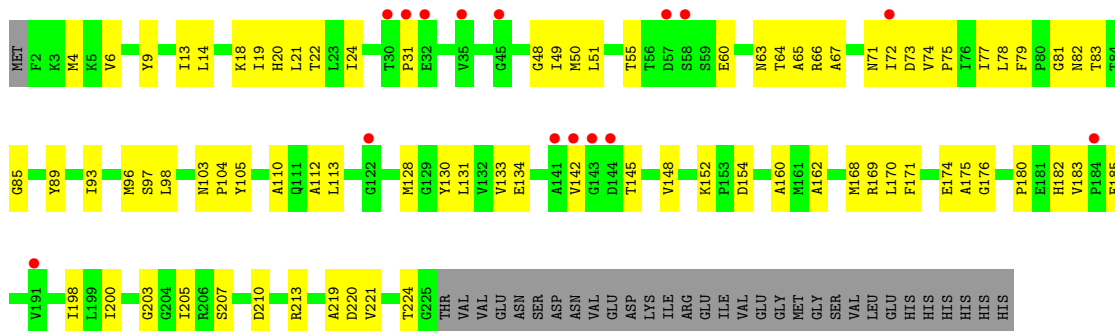


• Molecule 1: Geranylgeranylglyceryl phosphate synthase

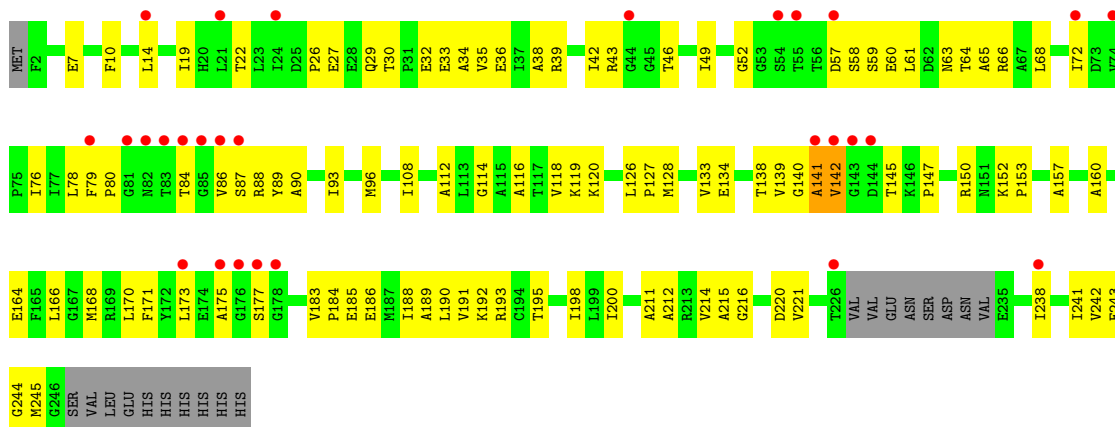




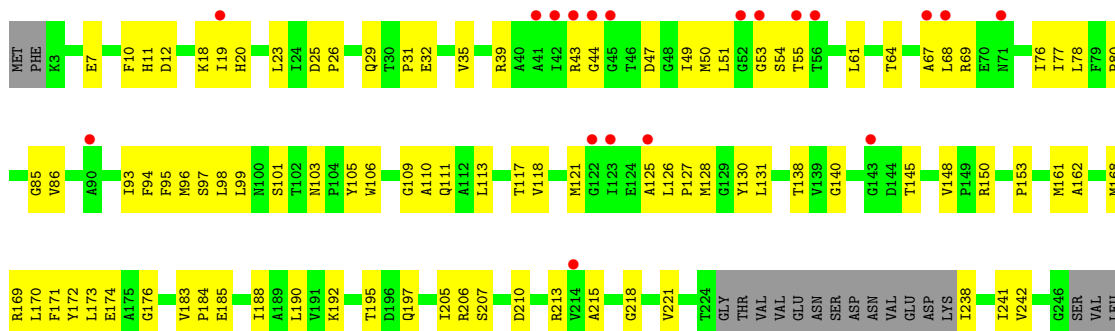
• Molecule 1: Geranylgeranylglyceryl phosphate synthase



• Molecule 1: Geranylgeranylglyceryl phosphate synthase



• Molecule 1: Geranylgeranylglyceryl phosphate synthase



GLU
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.20Å 62.78Å 136.26Å 90.00° 96.35° 90.00°	Depositor
Resolution (Å)	47.96 – 2.70 47.96 – 2.70	Depositor EDS
% Data completeness (in resolution range)	89.9 (47.96-2.70) 89.8 (47.96-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.232 , 0.321 0.234 , 0.324	Depositor DCC
R_{free} test set	1752 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	58.5	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 69.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9975	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.48	0/1756	0.68	0/2381
1	B	0.44	0/1617	0.70	0/2205
1	C	0.43	0/1653	0.65	0/2245
1	D	0.41	0/1638	0.65	0/2230
1	E	0.33	0/1734	0.58	0/2359
1	F	0.32	0/1692	0.56	0/2302
All	All	0.40	0/10090	0.64	0/13722

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1730	0	1717	61	0
1	B	1590	0	1537	71	0
1	C	1628	0	1596	44	0
1	D	1610	0	1575	69	0
1	E	1708	0	1654	75	0
1	F	1666	0	1637	83	0
2	A	8	0	0	0	0
2	B	6	0	0	1	0
2	C	20	0	0	3	0
2	D	5	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	2	0	0	2	0
2	F	2	0	0	0	0
All	All	9975	0	9716	389	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 20.

All (389) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:205:ILE:HD11	1:D:224:THR:HG22	1.27	1.15
1:F:98:LEU:HD21	1:F:106:TRP:HB3	1.45	0.95
1:D:170:LEU:HD22	1:D:198:ILE:HB	1.55	0.88
1:A:138:THR:HG22	1:B:152:LYS:HE2	1.58	0.85
1:C:26:PRO:HG2	1:C:61:LEU:HG	1.58	0.84
1:E:30:THR:HG22	1:E:33:GLU:HB2	1.60	0.84
1:D:205:ILE:HD11	1:D:224:THR:CG2	2.07	0.84
1:F:78:LEU:CB	1:F:93:ILE:HD12	2.08	0.83
1:B:120:LYS:HE3	1:B:121:MET:HE3	1.60	0.83
1:E:19:ILE:HB	1:E:221:VAL:HG12	1.60	0.83
1:C:164:GLU:OE1	2:C:301:HOH:O	1.99	0.81
1:D:79:PHE:CE1	1:D:96:MET:HE1	2.16	0.80
1:F:207:SER:HB2	1:F:210:ASP:H	1.46	0.79
1:D:182:HIS:HB3	1:D:205:ILE:HG22	1.63	0.79
1:B:39:ARG:O	1:B:39:ARG:NH1	2.17	0.77
1:D:20:HIS:HE1	1:D:22:THR:HG22	1.49	0.77
1:A:97:SER:HB3	1:A:168:MET:HE1	1.66	0.77
1:F:78:LEU:HB2	1:F:93:ILE:HD12	1.66	0.77
1:B:134:GLU:OE1	1:B:150:ARG:HG3	1.88	0.74
1:E:185:GLU:HA	1:E:188:ILE:HD12	1.70	0.72
1:B:39:ARG:HH12	1:B:43:ARG:N	1.88	0.72
1:E:59:SER:O	1:E:63:ASN:ND2	2.23	0.71
1:F:35:VAL:HG12	1:F:68:LEU:HD12	1.72	0.71
1:E:170:LEU:HD22	1:E:198:ILE:HB	1.73	0.71
1:D:20:HIS:CE1	1:D:22:THR:HG22	2.25	0.71
1:F:210:ASP:OD1	1:F:213:ARG:NH2	2.24	0.70
1:F:78:LEU:HB3	1:F:93:ILE:HD12	1.72	0.70
1:E:243:GLU:OE1	1:E:243:GLU:N	2.25	0.70
1:E:150:ARG:N	2:E:302:HOH:O	2.23	0.70
1:F:68:LEU:HB3	1:F:76:ILE:HD13	1.73	0.69
1:F:98:LEU:CD2	1:F:106:TRP:HB3	2.23	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:SER:O	1:B:99:LEU:HD12	1.91	0.69
1:A:207:SER:HB2	1:A:210:ASP:H	1.58	0.68
1:D:200:ILE:HG23	1:D:221:VAL:CG2	2.24	0.67
1:E:186:GLU:O	1:E:190:LEU:HG	1.93	0.67
1:F:19:ILE:HG23	1:F:221:VAL:HG22	1.76	0.67
1:A:30:THR:HG23	1:A:33:GLU:H	1.61	0.66
1:D:79:PHE:HE1	1:D:96:MET:HE1	1.60	0.66
1:E:26:PRO:HG2	1:E:61:LEU:HD12	1.77	0.66
1:F:51:LEU:HD22	1:F:76:ILE:HG22	1.78	0.66
1:F:80:PRO:HG3	1:F:93:ILE:HD11	1.79	0.65
1:A:98:LEU:HG	1:A:101:SER:HB2	1.78	0.65
1:F:64:THR:O	1:F:68:LEU:HD13	1.97	0.64
1:B:56:THR:OG1	1:B:57:ASP:N	2.28	0.64
1:E:211:ALA:HA	1:E:214:VAL:HG12	1.79	0.64
1:F:18:LYS:HD2	1:F:215:ALA:HB1	1.78	0.64
1:F:205:ILE:HD12	1:F:210:ASP:HB3	1.77	0.64
1:D:31:PRO:HB3	1:D:63:ASN:O	1.97	0.64
1:D:79:PHE:CD1	1:D:96:MET:HE1	2.33	0.64
1:C:79:PHE:CD1	1:C:96:MET:HE1	2.33	0.64
1:D:65:ALA:HB1	1:D:89:TYR:HB2	1.80	0.64
1:B:12:ASP:O	1:B:14:LEU:N	2.31	0.63
1:F:32:GLU:O	1:F:35:VAL:HG22	1.98	0.63
1:B:35:VAL:HA	1:B:68:LEU:HD21	1.80	0.63
1:B:182:HIS:ND1	1:B:183:VAL:O	2.31	0.62
1:E:78:LEU:HD11	1:E:87:SER:H	1.64	0.62
1:D:24:ILE:HG13	1:D:49:ILE:HD11	1.80	0.62
1:D:207:SER:HB2	1:D:210:ASP:H	1.63	0.62
1:F:127:PRO:HB2	1:F:168:MET:HA	1.81	0.62
1:B:30:THR:HG22	1:B:31:PRO:HD2	1.81	0.62
1:E:36:GLU:OE1	1:E:39:ARG:NH1	2.32	0.62
1:C:173:LEU:HD22	1:C:187:MET:HE3	1.81	0.62
1:E:108:ILE:HG12	1:F:99:LEU:HB3	1.82	0.62
1:F:195:THR:HG22	1:F:197:GLN:HG2	1.82	0.62
1:A:24:ILE:HG22	1:A:25:ASP:H	1.64	0.61
1:B:32:GLU:O	1:B:35:VAL:N	2.32	0.61
1:B:175:ALA:HB2	1:B:183:VAL:HG23	1.82	0.61
1:C:24:ILE:O	1:C:51:LEU:HA	2.00	0.61
1:F:140:GLY:O	1:F:145:THR:HG22	1.99	0.61
1:B:33:GLU:O	1:B:37:ILE:HG13	2.00	0.61
1:F:95:PHE:HE2	1:F:118:VAL:HG21	1.65	0.61
1:E:76:ILE:HD11	1:E:90:ALA:HA	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:192:LYS:NZ	1:F:218:GLY:O	2.24	0.61
1:C:23:LEU:HD12	1:C:50:MET:O	2.01	0.60
1:D:174:GLU:OE2	1:D:176:GLY:N	2.27	0.60
1:E:27:GLU:HA	1:E:60:GLU:OE2	2.01	0.60
1:F:54:SER:OG	1:F:55:THR:N	2.27	0.60
1:E:134:GLU:HB2	1:E:150:ARG:HG2	1.83	0.60
1:F:98:LEU:HG	1:F:101:SER:HB2	1.84	0.60
1:C:33:GLU:OE1	1:C:33:GLU:N	2.34	0.60
1:E:173:LEU:HD23	1:E:183:VAL:HG11	1.84	0.60
1:F:78:LEU:HB3	1:F:93:ILE:CD1	2.33	0.59
1:F:35:VAL:O	1:F:39:ARG:HG3	2.02	0.59
1:C:78:LEU:HB2	1:C:93:ILE:HG13	1.84	0.59
1:F:51:LEU:CD2	1:F:76:ILE:HG22	2.33	0.59
1:E:112:ALA:HA	1:F:162:ALA:HB2	1.85	0.59
1:E:145:THR:HG23	1:E:147:PRO:HD3	1.84	0.59
1:D:205:ILE:HB	1:D:210:ASP:HB3	1.85	0.58
1:F:185:GLU:HA	1:F:188:ILE:HD12	1.85	0.58
1:C:169:ARG:NE	2:C:304:HOH:O	2.28	0.58
1:B:42:ILE:HA	1:B:46:THR:HG22	1.84	0.58
1:E:157:ALA:HA	1:E:191:VAL:HG12	1.84	0.58
1:C:90:ALA:HB3	1:C:123:ILE:HD11	1.84	0.58
1:F:19:ILE:CD1	1:F:47:ASP:HB2	2.34	0.58
1:D:131:LEU:HD21	1:D:148:VAL:HG21	1.86	0.58
1:D:48:GLY:HA2	1:D:74:VAL:HB	1.87	0.57
1:D:19:ILE:HG12	1:D:221:VAL:HG12	1.86	0.57
1:E:150:ARG:NH1	1:E:184:PRO:HG3	2.19	0.57
1:D:98:LEU:HD13	1:D:130:TYR:HD1	1.69	0.57
1:D:6:VAL:HG11	1:D:75:PRO:HB3	1.85	0.57
1:E:68:LEU:HD23	1:E:72:ILE:HD11	1.85	0.57
1:F:43:ARG:HD3	1:F:44:GLY:N	2.19	0.57
1:B:33:GLU:HG2	1:B:37:ILE:HG13	1.86	0.57
1:B:33:GLU:O	1:B:37:ILE:N	2.24	0.57
1:F:31:PRO:HA	1:F:64:THR:HG23	1.86	0.57
1:B:38:ALA:O	1:B:42:ILE:HD12	2.05	0.57
1:D:78:LEU:HD11	1:D:93:ILE:HD12	1.87	0.57
1:F:78:LEU:CB	1:F:93:ILE:CD1	2.81	0.57
1:B:33:GLU:HA	1:B:36:GLU:HB2	1.88	0.56
1:B:134:GLU:OE1	1:B:150:ARG:CG	2.53	0.56
1:F:69:ARG:HE	1:F:69:ARG:C	2.13	0.56
1:F:150:ARG:CZ	1:F:184:PRO:HG3	2.35	0.56
1:C:183:VAL:HG13	1:C:187:MET:HE2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:153:PRO:HB2	1:E:190:LEU:HD12	1.87	0.56
1:A:67:ALA:O	1:A:71:ASN:ND2	2.32	0.56
1:B:13:ILE:HG22	1:B:17:ARG:H	1.69	0.56
1:C:134:GLU:HB3	1:C:150:ARG:HG3	1.88	0.56
1:C:195:THR:OG1	1:C:197:GLN:HG2	2.05	0.56
1:E:22:THR:OG1	1:E:49:ILE:HD13	2.06	0.55
1:B:166:LEU:HD22	1:D:168:MET:HE3	1.87	0.55
1:F:10:PHE:CE1	1:F:19:ILE:HG12	2.42	0.55
1:F:169:ARG:C	1:F:170:LEU:HD12	2.31	0.55
1:C:139:VAL:O	1:C:145:THR:HG21	2.06	0.55
1:C:211:ALA:HB2	1:C:241:ILE:HG23	1.87	0.55
1:E:200:ILE:HG23	1:E:221:VAL:HG23	1.89	0.55
1:B:13:ILE:HG23	1:B:16:GLU:HB3	1.88	0.55
1:F:171:PHE:HB2	1:F:197:GLN:HE21	1.72	0.55
1:A:48:GLY:HA2	1:A:74:VAL:CG1	2.36	0.55
1:B:13:ILE:HG22	1:B:17:ARG:N	2.21	0.55
1:D:60:GLU:N	1:D:60:GLU:OE1	2.40	0.55
1:F:10:PHE:HE1	1:F:19:ILE:HG12	1.72	0.55
1:F:78:LEU:HD13	1:F:86:VAL:HA	1.89	0.55
1:B:111:GLN:CD	1:B:111:GLN:H	2.14	0.54
1:F:171:PHE:HB2	1:F:197:GLN:NE2	2.22	0.54
1:A:134:GLU:HG2	1:A:149:PRO:HA	1.88	0.54
1:B:185:GLU:HB3	1:B:217:ALA:HB2	1.89	0.54
1:C:4:MET:HG2	1:C:5:LYS:N	2.23	0.54
1:C:42:ILE:HA	1:C:46:THR:HG22	1.88	0.54
1:F:23:LEU:HD12	1:F:50:MET:O	2.07	0.54
1:D:98:LEU:HD21	1:D:145:THR:HG22	1.90	0.54
1:E:14:LEU:HD23	1:E:14:LEU:O	2.07	0.54
1:C:96:MET:HA	1:C:128:MET:O	2.06	0.54
1:D:134:GLU:OE2	2:D:301:HOH:O	2.19	0.54
1:C:160:ALA:HA	1:C:171:PHE:CE1	2.42	0.54
1:D:154:ASP:HB2	2:D:304:HOH:O	2.07	0.53
1:A:179:ALA:O	1:A:206:ARG:NH2	2.41	0.53
1:F:10:PHE:C	1:F:12:ASP:H	2.17	0.53
1:D:31:PRO:HA	1:D:64:THR:HG22	1.90	0.53
1:E:38:ALA:O	1:E:42:ILE:HG23	2.08	0.53
1:F:26:PRO:HG2	1:F:61:LEU:HD12	1.91	0.53
1:F:78:LEU:HD22	1:F:85:GLY:O	2.08	0.53
1:A:74:VAL:HG13	1:A:75:PRO:HD2	1.91	0.53
1:E:39:ARG:HA	1:E:42:ILE:HG12	1.91	0.53
1:A:56:THR:OG1	1:A:57:ASP:N	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:ARG:HG3	1:A:89:TYR:CE1	2.44	0.53
1:A:145:THR:O	1:A:147:PRO:HD3	2.09	0.52
1:A:111:GLN:OE1	1:A:111:GLN:N	2.38	0.52
1:B:117:THR:O	1:B:120:LYS:HG3	2.09	0.52
1:E:138:THR:HG22	1:E:142:VAL:HG23	1.91	0.52
1:B:96:MET:HB3	1:B:128:MET:HB3	1.91	0.52
1:B:39:ARG:NH1	1:B:42:ILE:HB	2.25	0.52
1:E:80:PRO:HG3	1:E:93:ILE:HD11	1.91	0.52
1:E:116:ALA:O	1:E:120:LYS:HG2	2.09	0.52
1:E:166:LEU:HD22	1:F:168:MET:HE2	1.91	0.52
1:F:105:TYR:HA	1:F:109:GLY:HA3	1.92	0.52
1:A:135:PRO:HG2	1:A:180:PRO:HD2	1.93	0.51
1:E:32:GLU:HA	1:E:35:VAL:HG12	1.91	0.51
1:D:98:LEU:HD13	1:D:130:TYR:CD1	2.44	0.51
1:D:9:TYR:O	1:D:13:ILE:HG13	2.09	0.51
1:A:9:TYR:CZ	1:A:13:ILE:HD11	2.45	0.51
1:E:243:GLU:CD	1:E:244:GLY:N	2.69	0.51
1:F:131:LEU:HD21	1:F:148:VAL:HG21	1.92	0.51
1:A:130:TYR:OH	1:A:174:GLU:OE1	2.28	0.51
1:D:200:ILE:HG12	1:D:221:VAL:HG22	1.91	0.51
1:D:82:ASN:OD1	1:D:83:THR:N	2.39	0.51
1:A:134:GLU:HG3	1:A:150:ARG:HG3	1.93	0.51
1:A:239:ARG:HG2	1:A:243:GLU:OE2	2.10	0.51
1:B:50:MET:HA	1:B:77:ILE:O	2.10	0.50
1:B:88:ARG:NH1	1:B:89:TYR:OH	2.45	0.50
1:E:39:ARG:NH2	1:E:43:ARG:HH12	2.09	0.50
1:A:200:ILE:HG12	1:A:221:VAL:HB	1.92	0.50
1:F:19:ILE:CG2	1:F:221:VAL:HG22	2.41	0.50
1:C:183:VAL:N	2:C:306:HOH:O	2.33	0.50
1:D:200:ILE:HA	1:D:221:VAL:HG22	1.94	0.50
1:B:33:GLU:HG3	1:B:36:GLU:HB2	1.94	0.50
1:C:236:ASP:OD1	1:C:236:ASP:N	2.44	0.50
1:C:163:ALA:HA	1:C:166:LEU:HD12	1.94	0.50
1:F:19:ILE:HD12	1:F:20:HIS:H	1.77	0.50
1:B:68:LEU:O	1:B:72:ILE:HG12	2.12	0.49
1:D:200:ILE:HG23	1:D:221:VAL:HG23	1.94	0.49
1:E:189:ALA:O	1:E:193:ARG:HG3	2.12	0.49
1:F:195:THR:CG2	1:F:197:GLN:HG2	2.40	0.49
1:B:56:THR:HG22	1:F:113:LEU:O	2.11	0.49
1:F:98:LEU:HA	1:F:130:TYR:O	2.12	0.49
1:C:94:PHE:HB3	1:C:128:MET:HE2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:VAL:HG13	1:B:123:ILE:HG22	1.93	0.49
1:F:51:LEU:HD22	1:F:76:ILE:CG2	2.42	0.49
1:A:96:MET:HA	1:A:128:MET:O	2.13	0.49
1:E:27:GLU:OE2	2:E:301:HOH:O	2.20	0.49
1:A:55:THR:HG21	1:D:113:LEU:HD13	1.95	0.49
1:B:192:LYS:HB2	1:B:199:LEU:HD23	1.94	0.49
1:C:224:THR:HG21	1:C:241:ILE:HD13	1.95	0.49
1:E:96:MET:HA	1:E:128:MET:O	2.13	0.49
1:B:30:THR:CG2	1:B:31:PRO:HD2	2.43	0.48
1:B:163:ALA:HB1	1:B:168:MET:HB2	1.94	0.48
1:B:207:SER:H	1:B:210:ASP:HB2	1.78	0.48
1:A:55:THR:HG23	1:A:82:ASN:ND2	2.28	0.48
1:B:169:ARG:HG3	1:B:169:ARG:HH11	1.78	0.48
1:E:57:ASP:OD1	1:E:58:SER:N	2.46	0.48
1:E:30:THR:CG2	1:E:33:GLU:HB2	2.39	0.48
1:E:64:THR:O	1:E:68:LEU:HD12	2.14	0.48
1:A:160:ALA:HA	1:A:171:PHE:CE1	2.48	0.48
1:C:98:LEU:HD11	1:C:132:VAL:HG23	1.95	0.48
1:F:49:ILE:HB	1:F:76:ILE:HG23	1.96	0.48
1:A:96:MET:SD	1:A:128:MET:HE2	2.54	0.48
1:B:26:PRO:HB2	1:B:60:GLU:HG3	1.94	0.48
1:B:96:MET:HA	1:B:128:MET:O	2.13	0.48
1:D:55:THR:OG1	1:D:82:ASN:HB2	2.13	0.48
1:D:75:PRO:HB2	1:D:77:ILE:HD11	1.96	0.48
1:D:18:LYS:HD3	1:D:219:ALA:O	2.13	0.48
1:E:119:LYS:HD2	1:F:161:MET:SD	2.54	0.48
1:C:80:PRO:HD2	1:C:96:MET:HE3	1.96	0.48
1:E:160:ALA:HA	1:E:171:PHE:CE2	2.48	0.48
1:E:215:ALA:HB2	1:E:245:MET:HG2	1.96	0.48
1:C:121:MET:O	1:C:123:ILE:N	2.38	0.48
1:F:130:TYR:HD1	1:F:172:TYR:HB3	1.78	0.48
1:D:160:ALA:HA	1:D:171:PHE:CE1	2.49	0.47
1:F:174:GLU:HG3	1:F:176:GLY:H	1.78	0.47
1:B:98:LEU:HG	1:B:101:SER:HB2	1.95	0.47
1:A:65:ALA:HB1	1:A:89:TYR:HB2	1.96	0.47
1:D:185:GLU:HG2	1:D:213:ARG:HG3	1.95	0.47
1:F:96:MET:HA	1:F:128:MET:O	2.14	0.47
1:A:121:MET:C	1:A:123:ILE:H	2.23	0.47
1:D:4:MET:HE3	1:D:9:TYR:HB2	1.95	0.47
1:F:94:PHE:HB3	1:F:96:MET:HE3	1.95	0.47
1:A:17:ARG:HH22	1:A:47:ASP:CG	2.20	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ASN:OD1	1:A:84:THR:HG22	2.14	0.47
1:A:239:ARG:CG	1:A:243:GLU:OE2	2.63	0.47
1:E:127:PRO:HB2	1:E:168:MET:HA	1.96	0.47
1:E:212:ALA:O	1:E:216:GLY:N	2.44	0.47
1:B:78:LEU:HD13	1:B:86:VAL:HA	1.97	0.47
1:B:94:PHE:HE1	1:B:126:LEU:HD12	1.78	0.47
1:C:131:LEU:HD22	1:C:171:PHE:HZ	1.80	0.47
1:D:175:ALA:HB3	1:D:203:GLY:HA3	1.97	0.47
1:E:175:ALA:HB2	1:E:183:VAL:HG22	1.95	0.47
1:E:242:VAL:N	1:E:243:GLU:OE1	2.48	0.47
1:E:200:ILE:HG23	1:E:221:VAL:CG2	2.44	0.47
1:B:42:ILE:O	1:B:44:GLY:N	2.48	0.47
1:B:105:TYR:HA	1:B:109:GLY:HA3	1.97	0.47
1:B:164:GLU:OE2	2:B:301:HOH:O	2.21	0.47
1:A:128:MET:C	1:A:168:MET:HE2	2.40	0.46
1:D:24:ILE:CG1	1:D:49:ILE:HD11	2.43	0.46
1:E:42:ILE:HA	1:E:46:THR:HG22	1.97	0.46
1:A:211:ALA:HB2	1:A:241:ILE:HG23	1.98	0.46
1:E:80:PRO:HD2	1:E:96:MET:HE3	1.98	0.46
1:A:134:GLU:HA	1:A:135:PRO:C	2.40	0.46
1:B:31:PRO:HG3	1:B:63:ASN:ND2	2.30	0.46
1:F:69:ARG:O	1:F:69:ARG:NE	2.40	0.46
1:F:78:LEU:HB2	1:F:93:ILE:CD1	2.42	0.46
1:D:72:ILE:HG22	1:D:73:ASP:H	1.80	0.46
1:F:207:SER:N	1:F:210:ASP:HB2	2.31	0.46
1:F:19:ILE:HD12	1:F:20:HIS:N	2.30	0.46
1:F:173:LEU:HD23	1:F:183:VAL:HG11	1.98	0.46
1:B:170:LEU:HG	1:B:198:ILE:HB	1.98	0.46
1:A:37:ILE:HG21	1:A:226:THR:HG23	1.97	0.45
1:F:206:ARG:O	1:F:241:ILE:HD11	2.16	0.45
1:A:118:VAL:HG13	1:A:123:ILE:HG23	1.99	0.45
1:B:38:ALA:HB1	1:B:49:ILE:HD11	1.99	0.45
1:B:90:ALA:HB3	1:B:123:ILE:HD11	1.99	0.45
1:A:162:ALA:O	1:A:163:ALA:C	2.59	0.45
1:A:7:GLU:HG3	1:A:126:LEU:HD22	1.98	0.45
1:B:213:ARG:HH11	1:B:213:ARG:HG3	1.81	0.45
1:D:78:LEU:HD13	1:D:85:GLY:O	2.17	0.45
1:E:78:LEU:HD13	1:E:86:VAL:HA	1.97	0.45
1:D:105:TYR:CZ	1:D:110:ALA:HB2	2.51	0.45
1:C:133:VAL:CG2	1:C:187:MET:HE1	2.47	0.45
1:D:19:ILE:CG1	1:D:221:VAL:HG12	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:MET:HA	1:D:128:MET:O	2.17	0.45
1:D:133:VAL:HG11	1:D:183:VAL:HG13	1.98	0.45
1:C:75:PRO:HA	1:C:91:ASP:OD2	2.17	0.45
1:E:58:SER:OG	1:E:84:THR:HB	2.17	0.45
1:B:25:ASP:OD1	1:B:54:SER:HB2	2.16	0.44
1:C:30:THR:HG23	1:C:33:GLU:OE1	2.16	0.44
1:E:66:ARG:HH21	1:E:88:ARG:HH12	1.64	0.44
1:A:72:ILE:HD11	1:A:76:ILE:HD11	1.98	0.44
1:E:7:GLU:HB2	1:E:126:LEU:HD21	1.98	0.44
1:A:38:ALA:O	1:A:42:ILE:HD12	2.18	0.44
1:E:133:VAL:HG12	1:E:134:GLU:H	1.83	0.44
1:E:198:ILE:HA	1:E:220:ASP:OD2	2.17	0.44
1:F:25:ASP:O	1:F:29:GLN:NE2	2.51	0.44
1:F:105:TYR:CZ	1:F:110:ALA:HB2	2.53	0.44
1:D:60:GLU:O	1:D:64:THR:HG23	2.17	0.44
1:D:185:GLU:HG2	1:D:213:ARG:CG	2.47	0.44
1:D:182:HIS:HB3	1:D:205:ILE:CG2	2.42	0.44
1:F:7:GLU:HG3	1:F:126:LEU:HD22	2.00	0.44
1:F:19:ILE:HD11	1:F:47:ASP:HB2	1.99	0.44
1:A:48:GLY:HA2	1:A:74:VAL:HG13	1.99	0.44
1:A:50:MET:HE3	1:A:50:MET:HB3	1.85	0.44
1:B:27:GLU:HA	1:B:60:GLU:HG2	1.99	0.44
1:E:10:PHE:CD1	1:E:19:ILE:HD13	2.53	0.44
1:A:168:MET:HE3	1:C:166:LEU:HD22	1.98	0.44
1:B:213:ARG:HG3	1:B:213:ARG:NH1	2.33	0.44
1:D:152:LYS:HE2	1:F:138:THR:OG1	2.18	0.44
1:B:79:PHE:CE1	1:B:96:MET:HE1	2.53	0.43
1:C:133:VAL:HG22	1:C:187:MET:HE1	2.00	0.43
1:D:14:LEU:HD11	1:D:220:ASP:HB3	2.00	0.43
1:F:97:SER:HA	1:F:111:GLN:NE2	2.33	0.43
1:A:158:ALA:HB1	1:C:112:ALA:O	2.18	0.43
1:B:39:ARG:HD2	1:B:39:ARG:HA	1.62	0.43
1:A:27:GLU:OE2	1:A:54:SER:OG	2.29	0.43
1:A:131:LEU:HD11	1:A:148:VAL:HG21	2.00	0.43
1:B:30:THR:HG22	1:B:31:PRO:CD	2.47	0.43
1:B:111:GLN:H	1:B:111:GLN:NE2	2.17	0.43
1:E:29:GLN:HG3	1:E:34:ALA:HB2	1.99	0.43
1:F:31:PRO:O	1:F:35:VAL:HG13	2.18	0.43
1:C:50:MET:HA	1:C:77:ILE:O	2.19	0.43
1:D:72:ILE:HG22	1:D:73:ASP:N	2.33	0.43
1:F:207:SER:HB2	1:F:210:ASP:N	2.25	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:LEU:HD13	1:A:86:VAL:HA	2.01	0.43
1:A:129:GLY:N	1:A:168:MET:HE2	2.34	0.43
1:F:77:ILE:HG21	1:F:94:PHE:CZ	2.53	0.43
1:A:40:ALA:HB1	1:A:235:GLU:HG3	2.01	0.43
1:E:19:ILE:O	1:E:221:VAL:HA	2.19	0.43
1:E:140:GLY:HA2	1:E:145:THR:HG21	2.01	0.43
1:F:128:MET:HE3	1:F:128:MET:HB2	1.73	0.43
1:B:100:ASN:ND2	1:B:146:LYS:O	2.49	0.43
1:D:103:ASN:HD22	1:D:104:PRO:HD2	1.84	0.43
1:B:112:ALA:HA	1:D:162:ALA:HB2	2.01	0.42
1:A:237:LYS:HD3	1:A:237:LYS:HA	1.84	0.42
1:C:22:THR:HB	1:C:49:ILE:HD13	2.01	0.42
1:F:117:THR:O	1:F:121:MET:HG3	2.19	0.42
1:C:24:ILE:HD12	1:C:49:ILE:HD12	2.01	0.42
1:F:185:GLU:CD	1:F:185:GLU:H	2.28	0.42
1:D:182:HIS:CB	1:D:205:ILE:HG22	2.43	0.42
1:E:96:MET:HE2	1:E:96:MET:HB3	1.83	0.42
1:F:35:VAL:HG11	1:F:67:ALA:HB1	2.01	0.42
1:C:23:LEU:HA	1:C:50:MET:HG3	2.01	0.42
1:C:49:ILE:HG21	1:C:68:LEU:HD21	2.01	0.42
1:E:49:ILE:O	1:E:76:ILE:HA	2.19	0.42
1:E:114:GLY:O	1:E:118:VAL:HG12	2.20	0.42
1:E:152:LYS:HD2	1:E:152:LYS:HA	1.83	0.42
1:D:24:ILE:HB	1:D:51:LEU:HD12	2.01	0.42
1:E:241:ILE:C	1:E:243:GLU:OE1	2.62	0.42
1:F:153:PRO:HB2	1:F:190:LEU:HD13	2.01	0.42
1:A:96:MET:SD	1:A:128:MET:CE	3.08	0.42
1:A:113:LEU:HD21	1:D:142:VAL:HG13	2.02	0.42
1:D:169:ARG:C	1:D:170:LEU:HD23	2.44	0.42
1:E:139:VAL:C	1:E:141:ALA:H	2.27	0.42
1:F:238:ILE:O	1:F:242:VAL:HG23	2.19	0.42
1:E:238:ILE:O	1:E:242:VAL:HG23	2.20	0.42
1:B:9:TYR:O	1:B:12:ASP:HB2	2.20	0.41
1:B:158:ALA:HB1	1:D:112:ALA:O	2.20	0.41
1:C:4:MET:HE3	1:C:9:TYR:HB2	2.03	0.41
1:E:164:GLU:CD	1:E:195:THR:HB	2.45	0.41
1:B:6:VAL:O	1:B:9:TYR:HB3	2.20	0.41
1:C:195:THR:HG21	1:C:197:GLN:NE2	2.35	0.41
1:F:93:ILE:CG2	1:F:125:ALA:HA	2.51	0.41
1:E:52:GLY:HA3	1:E:79:PHE:HB3	2.02	0.41
1:A:153:PRO:HB2	1:A:190:LEU:HD23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:ILE:HG23	1:B:210:ASP:HB3	2.02	0.41
1:C:19:ILE:O	1:C:221:VAL:HA	2.20	0.41
1:A:103:ASN:HA	1:A:104:PRO:HD3	1.96	0.41
1:B:13:ILE:CG2	1:B:17:ARG:N	2.84	0.41
1:C:131:LEU:HD23	1:C:173:LEU:HD23	2.01	0.41
1:D:6:VAL:HA	1:D:9:TYR:HB3	2.02	0.41
1:D:97:SER:HB3	1:D:168:MET:HE1	2.01	0.41
1:D:128:MET:HE3	1:D:128:MET:HB2	1.72	0.41
1:E:26:PRO:HA	1:E:29:GLN:NE2	2.36	0.41
1:E:65:ALA:HB1	1:E:89:TYR:HB2	2.02	0.41
1:A:61:LEU:CD1	1:A:85:GLY:HA2	2.51	0.41
1:A:117:THR:O	1:A:121:MET:HG3	2.20	0.41
1:A:213:ARG:NE	1:A:213:ARG:HA	2.35	0.41
1:B:25:ASP:HB3	1:B:28:GLU:HB2	2.01	0.41
1:D:67:ALA:O	1:D:71:ASN:HB2	2.20	0.41
1:E:185:GLU:H	1:E:185:GLU:CD	2.29	0.41
1:F:103:ASN:OD1	1:F:103:ASN:C	2.63	0.41
1:B:24:ILE:N	1:B:24:ILE:HD12	2.36	0.41
1:E:192:LYS:HA	1:E:192:LYS:HD2	1.78	0.41
1:A:68:LEU:O	1:A:72:ILE:HG12	2.21	0.40
1:A:200:ILE:HG12	1:A:221:VAL:CG2	2.52	0.40
1:B:128:MET:HE3	1:B:128:MET:HB2	1.80	0.40
1:D:55:THR:OG1	1:D:81:GLY:O	2.39	0.40
1:B:152:LYS:HB3	1:B:152:LYS:HE3	1.82	0.40
1:D:21:LEU:HB3	1:D:50:MET:HG3	2.03	0.40
1:A:161:MET:O	1:A:162:ALA:C	2.64	0.40
1:E:138:THR:O	1:E:142:VAL:HG23	2.21	0.40
1:A:100:ASN:OD1	1:A:131:LEU:HD22	2.22	0.40
1:C:4:MET:HG2	1:C:5:LYS:H	1.86	0.40
1:D:50:MET:HE3	1:D:50:MET:HB3	1.93	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	227/256 (89%)	207 (91%)	16 (7%)	4 (2%)	6	18
1	B	219/256 (86%)	192 (88%)	23 (10%)	4 (2%)	6	18
1	C	218/256 (85%)	195 (89%)	16 (7%)	7 (3%)	3	7
1	D	222/256 (87%)	204 (92%)	16 (7%)	2 (1%)	14	35
1	E	233/256 (91%)	206 (88%)	24 (10%)	3 (1%)	9	25
1	F	227/256 (89%)	207 (91%)	18 (8%)	2 (1%)	14	35
All	All	1346/1536 (88%)	1211 (90%)	113 (8%)	22 (2%)	7	20

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	13	ILE
1	C	67	ALA
1	D	66	ARG
1	B	53	GLY
1	C	4	MET
1	C	88	ARG
1	E	142	VAL
1	E	177	SER
1	B	43	ARG
1	C	68	LEU
1	C	98	LEU
1	E	141	ALA
1	F	53	GLY
1	D	180	PRO
1	A	162	ALA
1	F	11	HIS
1	A	203	GLY
1	B	108	ILE
1	A	108	ILE
1	C	122	GLY
1	C	203	GLY
1	A	135	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar

resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	177/201 (88%)	177 (100%)	0	100	100
1	B	156/201 (78%)	156 (100%)	0	100	100
1	C	158/201 (79%)	158 (100%)	0	100	100
1	D	157/201 (78%)	157 (100%)	0	100	100
1	E	169/201 (84%)	169 (100%)	0	100	100
1	F	165/201 (82%)	165 (100%)	0	100	100
All	All	982/1206 (81%)	982 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	71	ASN
1	C	71	ASN
1	C	197	GLN
1	D	103	ASN
1	D	197	GLN
1	E	71	ASN
1	E	182	HIS
1	F	63	ASN
1	F	197	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	233/256 (91%)	0.11	7 (3%)	52	49	30, 50, 99, 146	0
1	B	221/256 (86%)	0.38	14 (6%)	26	23	34, 62, 111, 207	0
1	C	226/256 (88%)	0.31	8 (3%)	47	43	34, 56, 101, 116	0
1	D	224/256 (87%)	0.42	15 (6%)	24	21	35, 65, 119, 154	0
1	E	237/256 (92%)	0.78	28 (11%)	9	7	44, 84, 128, 152	0
1	F	231/256 (90%)	0.66	19 (8%)	17	15	45, 85, 129, 182	0
All	All	1372/1536 (89%)	0.45	91 (6%)	24	21	30, 67, 120, 207	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	45	GLY	6.0
1	F	41	ALA	5.8
1	B	46	THR	4.9
1	E	142	VAL	4.6
1	E	85	GLY	4.4
1	E	143	GLY	4.3
1	C	61	LEU	4.2
1	E	79	PHE	4.2
1	D	143	GLY	4.0
1	E	72	ILE	4.0
1	A	52	GLY	3.9
1	E	238	ILE	3.9
1	C	235	GLU	3.8
1	C	2	PHE	3.6
1	F	143	GLY	3.6
1	E	141	ALA	3.5
1	A	227	VAL	3.5
1	E	55	THR	3.5
1	E	82	ASN	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	86	VAL	3.3
1	D	32	GLU	3.2
1	D	141	ALA	3.2
1	D	45	GLY	3.1
1	E	54	SER	3.1
1	B	56	THR	3.0
1	E	87	SER	3.0
1	D	57	ASP	3.0
1	E	176	GLY	2.9
1	E	173	LEU	2.9
1	B	47	ASP	2.8
1	F	44	GLY	2.8
1	B	84	THR	2.8
1	F	53	GLY	2.8
1	F	52	GLY	2.8
1	D	191	VAL	2.7
1	E	175	ALA	2.7
1	A	26	PRO	2.7
1	C	86	VAL	2.7
1	D	142	VAL	2.7
1	B	122	GLY	2.7
1	C	31	PRO	2.7
1	B	44	GLY	2.6
1	B	67	ALA	2.6
1	F	71	ASN	2.6
1	A	235	GLU	2.6
1	F	67	ALA	2.6
1	D	30	THR	2.6
1	E	83	THR	2.6
1	A	56	THR	2.5
1	E	81	GLY	2.5
1	C	83	THR	2.5
1	E	84	THR	2.5
1	D	122	GLY	2.5
1	D	35	VAL	2.5
1	A	83	THR	2.5
1	A	84	THR	2.5
1	F	43	ARG	2.5
1	F	45	GLY	2.5
1	F	214	VAL	2.5
1	D	58	SER	2.4
1	E	14	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	55	THR	2.4
1	F	19	ILE	2.3
1	F	42	ILE	2.3
1	B	37	ILE	2.3
1	E	74	VAL	2.3
1	C	236	ASP	2.3
1	F	68	LEU	2.3
1	E	226	THR	2.3
1	F	90	ALA	2.2
1	D	184	PRO	2.2
1	F	55	THR	2.2
1	F	56	THR	2.2
1	E	178	GLY	2.2
1	B	83	THR	2.2
1	B	85	GLY	2.2
1	C	67	ALA	2.2
1	B	38	ALA	2.1
1	E	24	ILE	2.1
1	E	144	ASP	2.1
1	F	125	ALA	2.1
1	D	31	PRO	2.1
1	E	21	LEU	2.1
1	B	53	GLY	2.1
1	D	72	ILE	2.1
1	E	177	SER	2.0
1	F	122	GLY	2.0
1	D	144	ASP	2.0
1	E	57	ASP	2.0
1	E	44	GLY	2.0
1	F	123	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.